

UNITED STATES PATENT AND TRADEMARK OFFICE
CERTIFICATE OF CORRECTION

PATENT NO. : 7,235,576 B1
APPLICATION NO. : 10/042203
DATED : June 26, 2007
INVENTOR(S) : Riedl et al.

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Page 1 of 1

It is certified that error appears in the above-identified patent and that said Letters Patent is hereby corrected as shown below:

In column 3, line 31, reads "(CH₂)_m- and", should read -- (CH₂)_m-, and --.

In column 4, line 31, reads "methanesulphonic," should read -- methanesulfonic --.

In column 9, line 27, reads "ie.," should read -- i.e., --.

In column 9, line 39, reads "enclosure", should read -- disclosure --.

In column 17, line 1, reads "under stream", should read -- under a stream --.

In column 18, lines 39 and 40, reads "Aromatic", should read -- Aromatic --.

In column 20, line 29, reads "Methylcarbamoly", should read -- Methylcarbamoyl --.

In column 21, line 27, reads "(N-methylcarbamoly)", should read
-- (N-Methylcarbamoyl) --.

In column 21, line 38, reads "(N-methylcarbamoly)", should read
-- (N-Methylcarbamoyl) --.

In column 23, line 62, reads "then mixture", should read -- then the mixture --.

In column 30, line 4, reads "(4-(4-Methylsulfonylphenoxy))", should read
-- (4-(4-Methylsulfonylphenoxy) --.

In column 35, line 58, reads "heat", should read -- heated --.

In column 36, line 44, reads "ω-(Aroylamino)phenyl", should read
-- ω-(Arylamino)phenyl --.

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Page 2 of 2

It is certified that error appears in the above-identified patent and that said Letters Patent is hereby corrected as shown below:

In column 42, line 29, reads "(triflouromethyl)", should read -- trifluoromethyl --.

*ADD CORRECTIONS ON ATTACHED
PAGES

[REDACTED]

UNITED STATES PATENT AND TRADEMARK OFFICE

CERTIFICATE OF CORRECTION

PATENT NO.: 7,235,576 B1

Application No. 10/642,203

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DATED: June 26, 2007

INVENTOR(S): Riedl et al.

It is certified that error appears in the above-identified patent and that said Letters Patent are hereby corrected as shown below:

In columns 92-98, replace claims 2-47 with the claims 2-17 as follows:

2. A pharmaceutically acceptable salt of claim 1, which is

a) a basic salt of an organic acid or an inorganic acid which is hydrochloric acid, hydrobromic acid, sulfuric acid, phosphoric acid, methanesulfonic acid, trifluoromethanesulfonic acid, benzenesulfonic acid, p-toluene sulfonic acid (tosylate salt), 1-naphthalene sulfonic acid, 2-naphthalene sulfonic acid, acetic acid, trifluoroacetic acid, malic acid, tartaric acid, citric acid, lactic acid, oxalic acid, succinic acid, fumaric acid, maleic acid, benzoic acid, salicylic acid, phenylacetic acid, or mandelic acid; or

b) an acid salt of an organic or inorganic base containing an alkali metal cation, an alkaline earth metal cation, an ammonium cation, an aliphatic substituted ammonium cation or an aromatic substituted ammonium cation.

3. A pharmaceutically acceptable salt of claim 1 which is a basic salt of an organic acid or an inorganic acid which is hydrochloric acid, hydrobromic acid, sulfuric acid, phosphoric acid, methanesulfonic acid, trifluoromethanesulfonic acid, benzenesulfonic acid, p-toluene sulfonic acid (tosylate salt), 1-naphthalene sulfonic acid, 2-naphthalene sulfonic acid, acetic acid, trifluoroacetic acid, malic acid, tartaric acid, citric acid, lactic acid, oxalic acid, succinic acid, fumaric acid, maleic acid, benzoic acid, salicylic acid, phenylacetic acid, or mandelic acid.

4. A pharmaceutically acceptable salt which is the tosylate salt of

N-(5-*tert*-butyl-2-methoxy phenyl)-*N'*-(4-(4-methoxy-3-(*N*-methylcarbamoyl) phenoxy) phenyl) urea,
N-(2-methoxy-5-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl)

urea,

N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-carbamoyl-4-pyridyloxy) phenyl) urea,

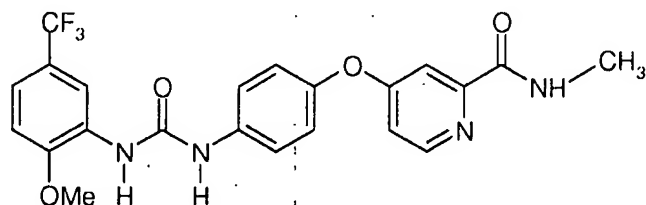
N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl)
urea; or

N-(2-methoxy-4-chloro-5-(trifluoromethyl) phenyl)-*N'*-(3-(2-(*N*-methylcarbamoyl)-4-pyridyloxy)

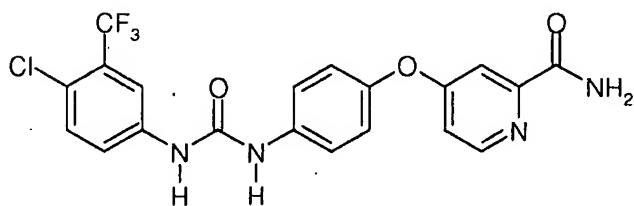
phenyl) urea.

5. A pharmaceutically acceptable salt of a compound which is:

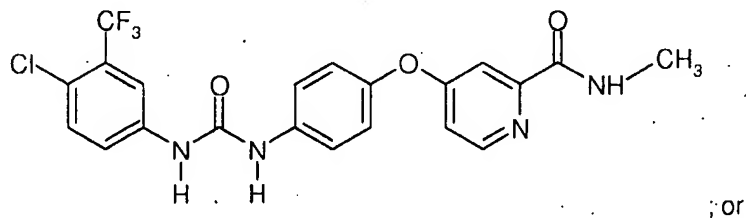
N-(2-methoxy-5-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:



N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-carbamoyl-4-pyridyloxy) phenyl) urea of the formula:

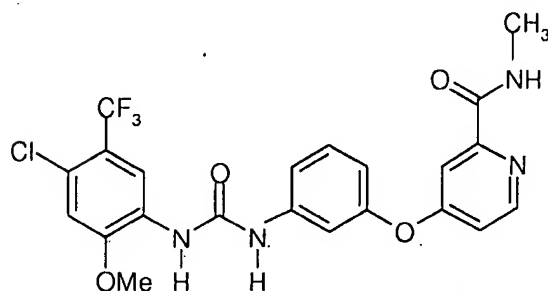


N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:



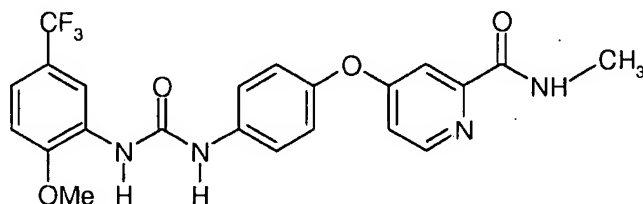
; or

N-(2-methoxy-4-chloro-5-(trifluoromethyl) phenyl)-*N'*-(3-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:

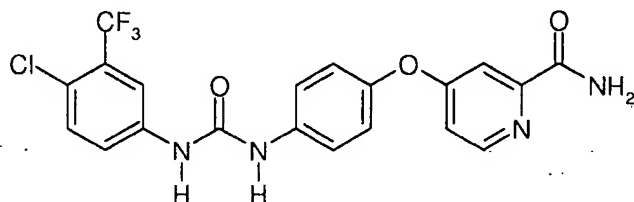


6. A pharmaceutically acceptable salt of claim 5 which is a basic salt of an organic acid or an inorganic acid which is hydrochloric acid, hydrobromic acid, sulfuric acid, phosphoric acid, methanesulfonic acid, trifluoromethanesulfonic acid, benzenesulfonic acid, p-toluene sulfonic acid (tosylate salt), 1-naphthalene sulfonic acid, 2-naphthalene sulfonic acid, acetic acid, trifluoroacetic acid, malic acid, tartaric acid, citric acid, lactic acid, oxalic acid, succinic acid, fumaric acid, maleic acid, benzoic acid, salicylic acid, phenylacetic acid, or mandelic acid.

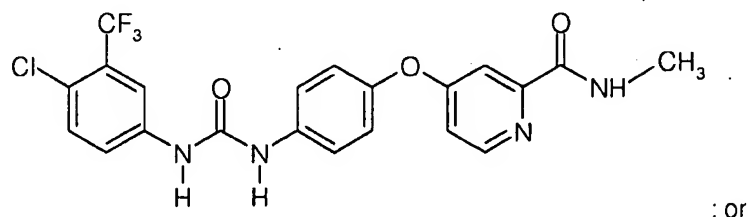
7. A pharmaceutically acceptable salt which is the tosylate salt of
N-(2-methoxy-5-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:



N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-carbamoyl-4-pyridyloxy) phenyl) urea of the formula:



N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:

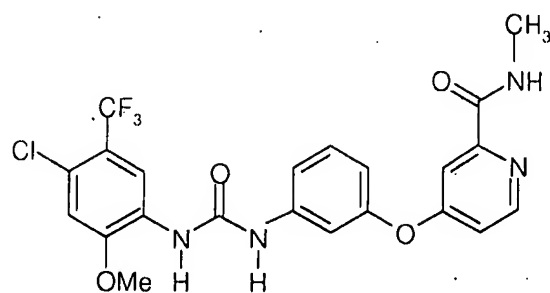


; or

N-(2-methoxy-4-chloro-5-(trifluoromethyl) phenyl)-*N'*-(3-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:

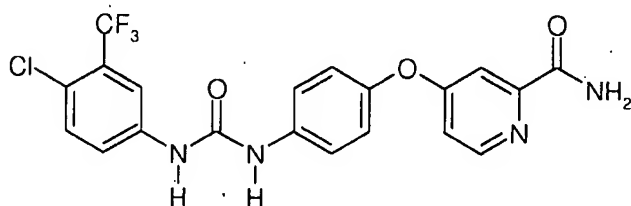
Nok

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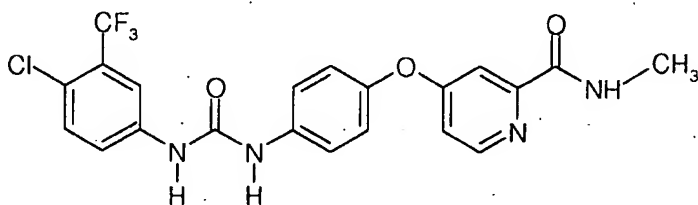


8. A pharmaceutically acceptable salt of a compound which is:

N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-carbamoyl-4-pyridyloxy) phenyl) urea of the formula:

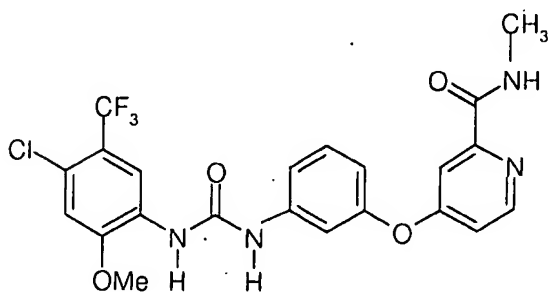


N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:



; or

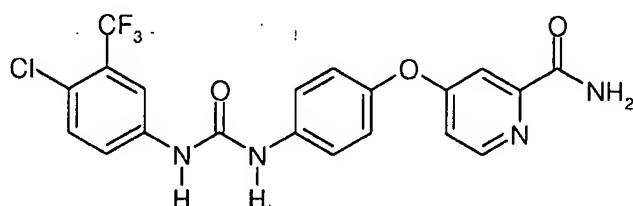
N-(2-methoxy-4-chloro-5-(trifluoromethyl) phenyl)-*N'*-(3-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:



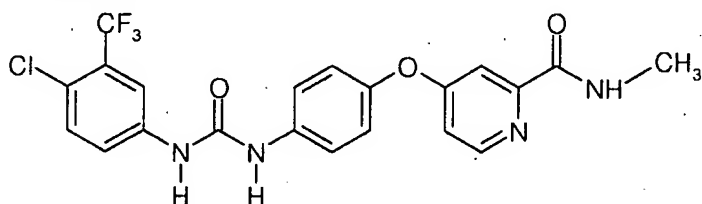
9. A pharmaceutically acceptable salt of claim 8 which is a basic salt of an organic acid or an inorganic acid which is hydrochloric acid, hydrobromic acid, sulfuric acid, phosphoric acid, methanesulfonic acid, trifluoromethanesulfonic acid, benzenesulfonic acid, p-toluene sulfonic acid (tosylate salt), 1-naphthalene sulfonic acid, 2-naphthalene sulfonic acid, acetic acid, trifluoroacetic acid, malic acid, tartaric acid, citric acid, lactic acid, oxalic acid, succinic acid, fumaric acid, maleic acid, benzoic acid, salicylic acid, phenylacetic acid, or mandelic acid.

10. A pharmaceutically acceptable salt which is the tosylate salt of

N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-carbamoyl-4-pyridyloxy) phenyl) urea of the formula:

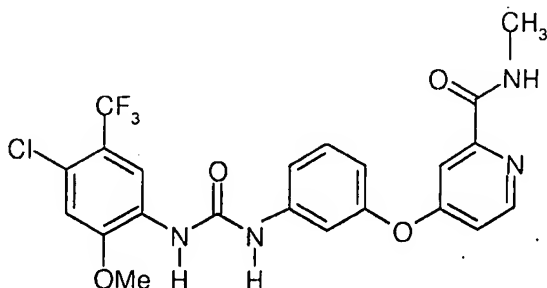


N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:



; or

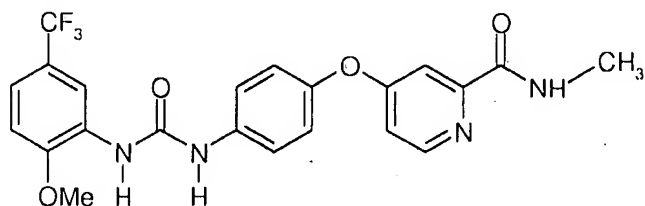
N-(2-methoxy-4-chloro-5-(trifluoromethyl) phenyl)-*N'*-(3-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:



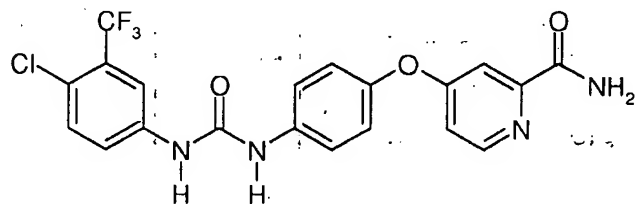
11. A pharmaceutically acceptable salt of a compound which is:

N-(2-methoxy-5-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:

Note

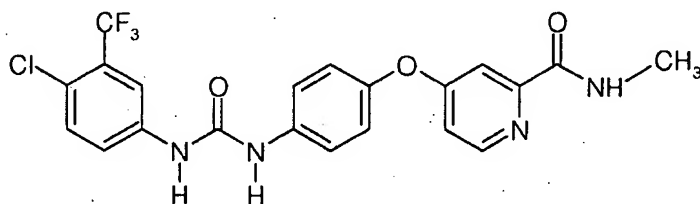


N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-carbamoyl-4-pyridyloxy) phenyl) urea of the formula:



, or

N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:

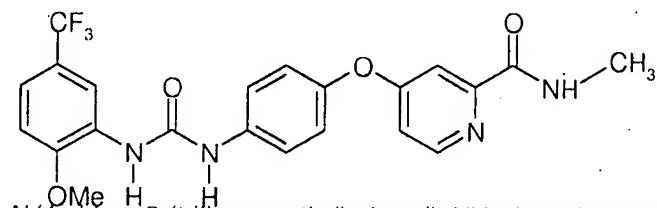


12. A pharmaceutically acceptable salt of claim 11 which is a basic salt of an organic acid or an inorganic acid which is hydrochloric acid, hydrobromic acid, sulfuric acid, phosphoric acid, methanesulfonic acid, trifluoromethanesulfonic acid, benzenesulfonic acid, p-toluene sulfonic acid (tosylate salt), 1-naphthalene sulfonic acid, 2-naphthalene sulfonic acid, acetic acid, trifluoroacetic acid, malic acid, tartaric acid, citric acid, lactic acid, oxalic acid, succinic acid, fumaric acid, maleic acid, benzoic acid, salicylic acid, phenylacetic acid, or mandelic acid.

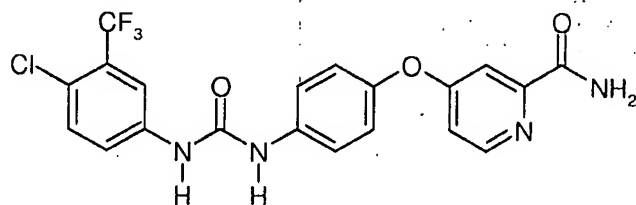
13. A pharmaceutically acceptable salt which is the tosylate salt of

N-(2-methoxy-5-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:

not

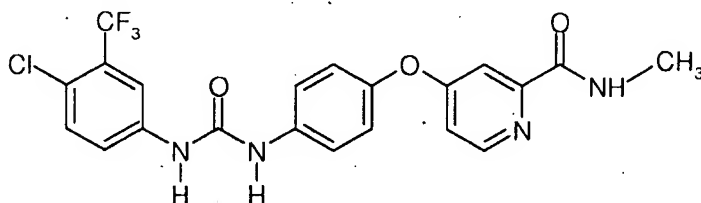


N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-carbamoyl-4-pyridyloxy) phenyl) urea of the formula:



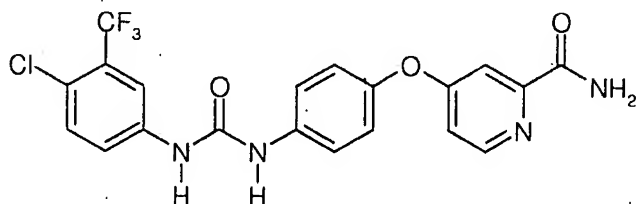
, or

N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula



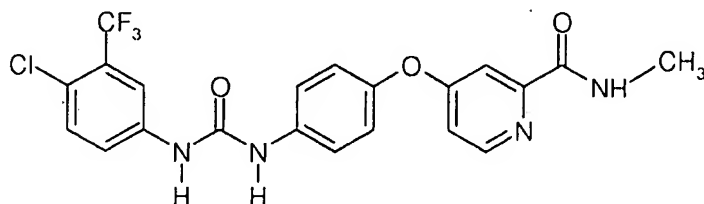
14. A pharmaceutically acceptable salt of a compound which is:

N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-carbamoyl-4-pyridyloxy) phenyl) urea of the formula:



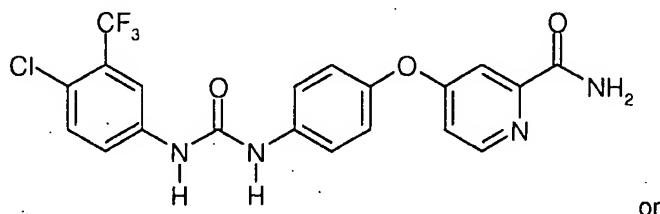
or

N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:

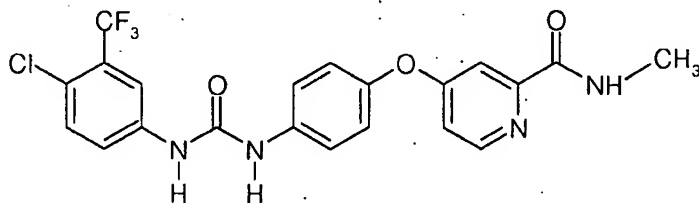


15. A pharmaceutically acceptable salt of claim 14 which is a basic salt of an organic acid or an inorganic acid which is hydrochloric acid, hydrobromic acid, sulfuric acid, phosphoric acid, methanesulfonic acid, trifluoromethanesulfonic acid, benzenesulfonic acid, p-toluene sulfonic acid (tosylate salt), 1-naphthalene sulfonic acid, 2-naphthalene sulfonic acid, acetic acid, trifluoroacetic acid, malic acid, tartaric acid, citric acid, lactic acid, oxalic acid, succinic acid, fumaric acid, maleic acid, benzoic acid, salicylic acid, phenylacetic acid, or mandelic acid.

16. A pharmaceutically acceptable salt which is the tosylate salt of
N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-carbamoyl-4-pyridyloxy) phenyl) urea of the formula:



N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula



17. A pharmaceutical composition comprising a pharmaceutically acceptable salt of claim 1, 2, 3, 4, 5, 6, 7, 8, 9, 10, 11, 12, 13, 14, 15 or 16 and pharmaceutically acceptable carrier.